PATENT NO. : 7,071,205 B2 APPLICATION NO. : 10/684229 Page 1 of 10

APPLICATION NO.: 10/684229
DATED: July 4, 2006
INVENTOR(S): Lin Zhi et al.

It is certified that error appears in the above-identified patent and that said Letters Patent is hereby corrected as shown below:

IN THE TITLE PAGES:

In Item [56] References Cited, in U.S. PATENT DOCUMENTS: please add the following reference: --6,001,846 A 12/1999 Edwards et al. 514/285--in 6,566,358 please replace "Zhi et al." with --Zhang et a1.--in 6,566,372 please replace "West et al." with --Zhi et al.--

In Item [56] References Cited, in OTHER PUBLICATIONS: in Hamann et al., please replace "dihyrdo" with --dihydro--

At column 8, Table A, row R¹, please replace "C₁-C haloalkyl" with --C₁-C₄ haloalkyl-at column 9, Table A, row R⁹, please replace "CONR^HR¹²" with --CONR¹¹R¹²-- at column 11, Table A, below row R¹⁶, please replace "R¹⁵" with --R¹⁶--

Please replace Claims 12, 13, 14, and 15 with the following Claims:

Col. 40

12. A compound of the formula:

(I)

wherein:

 R^1 is selected from the group of hydrogen, C_1 – C_4 alkyl, C_1 – C_4 haloalkyl, C_1 – C_4 heteroalkyl, COR^{11} , CO_2R^{11} , SO_2R^{11} , and $CONR^{11}R^{12}$;

 R^2 and R^3 each independently is selected from the group of hydrogen, C_1 – C_6 alkyl, and C_1 – C_6 haloalkyl; or

R² and R³ taken together form a cycloalkyl ring of from three to twelve carbons;

 R^4 through R^7 each independently is selected from the group of hydrogen, F, C1, Br, CN, QR^{11} , Q_1 – Q_4 alkyl, Q_1 – Q_4 haloalkyl, and Q_1 – Q_4 heteroalkyl; or

R⁵ and R⁷ taken together form a bond; or

R⁶ and R⁷ taken together are selected from the group of methylidene, mono-substituted methylidene, di-substituted methylidene and carbonyl;

PATENT NO. : 7.071,205 B2 **APPLICATION NO. : 10/684229**

Page 2 of 10

DATED : July 4, 2006 INVENTOR(S) : Lin Zhi et al.

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 R^8 through R^{10} each independently is selected from the group of hydrogen, F, Cl, Br, I, NO₂, CN, OR^{11} , $NR^{11}R^{12}$, SR^{11} , COR^{11} , CO_2R^{11} , $CONR^{11}R^{12}$, C_1 – C_8 alkyl, C_1 – C_8 heteroalkyl, C_1 – C_8 haloalkyl, allyl, C_2 – C_8 alkenyl and C_2 – C_8 alkynyl;

 R^{11} and R^{12} each is independently selected from the group of hydrogen, C_1 – C_4 alkyl, C₁-C₄ heteroalkyl, and C₁-C₄ haloalkyl;

R¹³ is hydrogen;

R¹⁴ and R¹⁶ taken together form a bond or "-O-" bridge;

R¹⁵, R¹⁷, R¹⁸, R¹⁹, R²⁰ each independently is selected from the group of hydrogen, F, C1, C_1 – C_4 alkyl, and C_1 – C_4 haloalkyl;

R²¹ is hydrogen; and n is 0, 1, 2, or 3;

or a pharmaceutically acceptable salt thereof.

Col. 41

13. A compound of the formula:

(I)

PATENT NO. : 7.071,205 B2 APPLICATION NO.: 10/684229

: July 4, 2006

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Page 3 of 10

wherein:

R¹ is selected from the group of hydrogen, C₁-C₄ alkyl, C₁-C₄ haloalkyl, C₁-C₄ heteroalkyl, COR¹¹, CO₂R¹¹, SO₂R¹¹, and CONR¹¹R¹²;

R² and R³ each independently is selected from the group of hydrogen, C₁-C₆ alkyl, and C₁-C₆ haloalkyl; or

R² and R³ taken together form a cycloalkyl ring of from three to twelve carbons;

R⁴ through R⁷ each independently is selected from the group of hydrogen, F, Cl, Br, CN, OR^{11} , C_1 - C_4 alkyl, C_1 - C_4 haloalkyl, and C_1 - C_4 heteroalkyl; or

R⁵ and R⁷ taken together form a bond; or

R⁶ and R⁷ taken together are selected from the group of methylidene, mono-substituted methylidene, di-substituted methylidene and carbonyl;

R⁸ through R¹⁰ each independently is selected from the group of hydrogen, F, Cl, Br, I, NO₂, CN, OR¹¹, NR¹¹R¹², SR¹¹, COR¹¹, CO₂R¹¹, CONR¹¹R¹², C₁-C₈ alkyl, C₁-C₈ heteroalkyl, C₁-C₈ haloalkyl, allyl, C₂-C₈ alkenyl and C₂-C₈ alkynyl;

 R^{11} and R^{12} each is independently selected from the group of hydrogen, C_1 – C_4 alkyl, C_1 – C_4 heteroalkyl, and C_1 – C_4 haloalkyl;

R¹³ is hydrogen;

R¹⁴, R¹⁵, R¹⁸, R¹⁹, R²⁰ each independently is selected from the group of hydrogen, F, Cl, C_1 – C_4 alkyl, and C_1 – C_4 haloalkyl;

R¹⁶ and R¹⁷ taken together are selected from the group of methylidene, mono-substituted methylidene, and di-substituted methylidene;

R²¹ is hydrogen; or

R²¹ and R²⁰ taken together form a bond;

n is 0, 1, 2, or 3;

or a pharmaceutically acceptable salt thereof.

PATENT NO.

: 7,071,205 B2

Page 4 of 10

APPLICATION NO. : 10/684229 DATED

: July 4, 2006

INVENTOR(S)

: Lin Zhi et al.

It is certified that error appears in the above-identified patent and that said Letters Patent is hereby corrected as shown below:

Col. 42

14. A compound of the formula:

(II)

wherein:

R1 is selected from the group of hydrogen, C1-C4 alkyl, C1-C4 haloalkyl, C1-C4 heteroalkyl, COR¹¹, CO₂R¹¹, SO₂R¹¹, and CONR¹¹R¹²;

R² and R³ each independently is selected from the group of hydrogen, C₁-C₆ alkyl, and C₁-C₆ haloalkyl; or

R² and R³ taken together form a cycloalkyl ring of from three to twelve carbons;

R⁴ through R⁷ each independently is selected from the group of hydrogen, F, Cl, Br, CN, OR^{11} , C_1 – C_4 alkyl, C_1 – C_4 haloalkyl, and C_1 – C_4 heteroalkyl; or

R⁵ and R⁷ taken together form a bond; or

R⁶ and R⁷ taken together are selected from the group of methylidene, mono-substituted methylidene, di-substituted methylidene and carbonyl;

 R^8 through R^{10} each independently is selected from the group of hydrogen, F, Cl, Br, I, NO₂, CN, OR¹¹, NR¹¹R¹², SR¹¹, COR¹¹, CO₂R¹¹, CONR¹¹R¹², C₁-C₈ alkyl, C₁-C₈ heteroalkyl, C1-C8 haloalkyl, allyl, C2-C8 alkenyl and C2-C8 alkynyl;

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: 7,071205 B2

APPLICATION NO.: 10/684229

DATED INVENTOR(S) : July 4, 2006 : Lin Zhi et al.

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R¹¹ and R¹² each is independently selected from the group of hydrogen, C₁-C₄ alkyl, C₁-C₄ heteroalkyl, and C₁-C₄ haloalkyl:

Page 5 of 10

R¹³ is hydrogen:

R¹⁴, R¹⁵, R¹⁷, R²⁰ each independently is selected from the group of hydrogen, F, Cl, C₁- C_4 alkyl, and C_1 – C_4 haloalkyl;

 R^{16} and R^{18} taken together form a bond when n is 1; R^{16} and R^{19} taken together form a bond when n is 0;

R²¹ is hydrogen; and

n is 0, 1, 2, or 3;

or a pharmaceutically acceptable salt thereof.

Col. 42-45

15. A Compound selected from the group of:

(±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-

chromeno[3,4-f]quinoline (compound 24);

 (\pm) -(51,1'u)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 25);

(+)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 27):

(-)-(51,1'l)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 28);

(±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-hydroxy-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 29);

PATENT NO. : 7.071,205 B2 APPLICATION NO.: 10/684229 DATED

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Page 6 of 10

- (±)-(51,1'u)-5-(3-methyl-2-cyclohexenyl)-9-hydroxy-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 30);
- (+)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-hydroxy-1,2-dihydro-2,2,4-trimethyl-5Hcbromeno[3,4-fjquinoline (compound 32);
- (-)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-hydroxy-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-fJquinoline (compound 33);
- (±)-(51,1'1)-5-3-methy1-2-cyclohexeny1)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 34);
- (±)-(51,1'u)-5-(3-methyl-2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 35);
- (+)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 37);
- (-)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 38);
- (±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-methoxy-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 39);
- (±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2-dimethyl-5Hchromeno[3,4-f]quinoline (compound 41):
- (\pm) -(51,1'u)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2-dimethyl-5Hchromeno[3,4-f]quinoline (compound 42);
- (±)-(51,1'1)-5-(3-methy1-2-cyclopenteny1)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-. chromeno[3,4-f]quinoline (compound 44);
- (±)-(51,1'u)-5-(3-methyl-2-cyclopentenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethy-5Hchromeno[3,4-f]quinoline (compound 45);

PATENT NO. : 7,071,205 B2 APPLICATION NO.: 10/684229

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(±)-(51,1'1)-5-(3,5,5-trimethyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 47);

Page 7 of 10

- (±)-(51,1'u)-5-(3,5,5-trimethyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 48);
- (±)-(51,1'1)-5-(3-methyl-2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 50);
- (±)-(51,1'u)-5-(3-methyl-2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 51);
- (±)-5-(3-methyl-3-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 52);
- (±)-5-(2-cyclopenta-1,3-dienyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4f]quinoline (compound 53);
- (±)-(51,1'1)-5-(3-ethy1-2-cyclohexeny1)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 55);
- (±)-(51,1'u)-5-(3-ethyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 56);
- (±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-7-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 58);
- (±)-(51,1'u)-5-(3-methy1-2-cyclohexeny1)-7-fluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 59);
- (±)-(51,1'1)-5-(3-ethyl-2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 61);
- (±)-(51,1'1)-5-(3-ethylidenecyclohexyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 62);

PATENT NO. : 7,071,205 B2 **APPLICATION NO. : 10/684229**.

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Page 8 of 10

- (±)-(51,1'1)-5-(3-methyl-3-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 63);
- (±)-(51,1'1)-5-(3-methyl-2-cyclobexenyl)-9-fluoro-1,2-dihydro-8-methoxy-2,2,4trimethyl-5H-chromeno[3,4-f]quinoline (compound 64);
- (±)-(51,1'u)-5-(3-methyl-2-cyclohexenyl)-9-fluoro-1,2-dihydro-8-methoxy-2,2,4trimethyl-5H-chromeno[3,4-f]quinoline (compound 65);
- (±)-(51,1'1)-5-(2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 67);
- (±)-(51,1'u)-5-(2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 68);
- (±)-5-(1-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4f]quinoline (compound 69);
- (±)-(51,1'1)-5-(2,3-dimethyl-2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 71): cy do penteny 1
- (+)-(51,1'1)-5-(2,3-dimethyl-2-myclopycolol)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 73);
- (-)-(51,1'1)-5-(2,3-dimethyl-2-cyclopentenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 74):
- (±)-(51,1'1)-5-(2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 75);
- (±)-(51,1'u)-5-(2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5Hchromeno[3,4-f]quinoline (compound 76);
- (±)-(51,1'1)-5-(2-cyclohexenyl)-7,9-difluoro-1,2,3,4-tetrahydro-2,2-dimethyl-4methylidene-5H-chromeno[3,4-f]quinoline (compound 77);

PATENT NO. : 7,071,205 B2 APPLICATION NO. : 10/684229

Page 9 of 10

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- (±)-(51,1'1)-5-(2-methylidenecyclohexyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 79):
- (±)-(51,1'u)-5-(2-methylidenecyclohexyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 80);
- (±)-(51,1'1)-5-(2-oxocyclohexyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 81);
- (±)-(51,1'u)-5-(2-oxocyclohexyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 82);
- (±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-9-methoxy-1,2-dihydro-1,2,2,4-tetramethyl-5H-chromeno[3,4-f]quinoline (compound 83);
- (±)-5-(2-cyclohexenyl)-9-fluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]-quinoline (compound 84);
- (±)-(51,1'l)-5-(2,3-dimethyl-2-cyclohexenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 85);
- (±)-5-(3-methylidene-cyclohexyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 87);
- (±)-(51,1'u)-5-(3-ethylidenecyclohexyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (compound 88);
- (±)-(51,1'1)-5-(2-cycloheptenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (Compound 89);
- (±)-(51,1'1)-5-(2-cyclooctenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (Compound 91);
- (±)-(51,1'u)-5-(2-cyclooctenyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (Compound 92);

PATENT NO.

: 7,071,205 B2

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(±)-(51,1'1)-5-(2,3-epoxy-3-methylcyclohexyl)-7,9-difluoro-1,2-dihydro-2,2,4-trimethyl-5H-chromeno[3,4-f]quinoline (Compound 94);

methylene-5H-chromeno[3,4-f]quinolin-3-ol (Compound 95);

(±)-(51,1'1)-5-(2,3-epoxy-2,3-dimethylcyclopentyl)-7,9-difluoro-1,2-dihydro-2,2,4trimethyl-5H-chromeno[3,4-f]quinoline (Compound 96);

(±)-(51,1'u)-5-(2,3-epoxy-3-methy1cyc1ohexy1)-7,9-dif1uoro-1,2-dihydro-2,2,4trimethy1-5H-chromeno[3,4-f]quinoline (Compound 97); and

(±)-(51,1'1)-5-(3-methyl-2-cyclohexenyl)-7,9-difluoro-1,2,3,4-tetrahydro-2,2-dimethyl-5H-chromeno[3,4-f]quinolin-4-one (Compound 98).

This certificate supersedes CANNAMA of Correction issued November 28, 2006.

April 3,2007. and

(attached and capted and applications)

Page 10 of 10

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